Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090815\
Data File : BE090747.D
Acq On : 8 Sep 2015 9:48
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 08 23:03:45 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	129	0.00
2		1,4-Dioxane	0.814	0.709	12.9	116	0.00
3		n-Nitrosodimethylamine	1.033	0.900	12.9	119	0.00
4	S	2-Fluorophenol	1.039	1.003	3.5	133	0.00
5	S	Phenol-d6	1.392	1.580	-13.5	154#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	143	0.00
7	S	Nitrobenzene-d5	0.330	0.344	-4.2	149	0.00
8		Nitrobenzene	0.365	0.379	-3.8	155#	0.00
9		Naphthalene	1.091	1.091	0.0	144	0.00
10		Hexachlorobutadiene	0.171	0.160	6.4	130	0.00
11		2-Methylnaphthalene	0.590	0.671	-13.7	168#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	161#	0.00
13	S	2,4,6-Tribromophenol	0.143	0.144	-0.7	191#	0.00
14	S	2-Fluorobiphenyl	1.385	1.366	1.4	161#	0.00
15		Acenaphthylene	7.992	8.430	-5.5	177#	0.00
16		Acenaphthene	1.335	1.344	-0.7	160#	0.00
17		Fluorene	1.666	1.707	-2.5	166#	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	153#	0.00
19		4-Bromophenyl-phenylether	0.170	0.180	-5.9	168#	0.00
20		Hexachlorobenzene	0.928	0.889	4.2	149	0.00
21		Pentachlorophenol	0.054	0.053	1.9	173#	0.00
22		Phenanthrene	1.237	1.283	-3.7	159#	-0.02
23		Anthracene	1.035	1.157	-11.8	178#	0.00
24		Fluoranthene	1.304	1.436	-10.1	182#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	159#	0.00
26		Pyrene	1.035	1.202	-16.1	194#	0.00
27	S	Terphenyl-d14	0.553	0.613	-10.8	184#	0.00
28		Benzo(a)anthracene	1.156	1.216	-5.2	176#	0.00
29		Chrysene	1.279	1.337	-4.5	168#	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.337	2.3	211#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.387	-17.4	202#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	173#	0.00
33		Benzo(b)fluoranthene	1.084	1.128	-4.1	185#	0.00
34		Benzo(k)fluoranthene	1.263	1.349	-6.8	187#	0.00
35	C	Benzo(a)pyrene	0.970	1.081	-11.4	206#	0.00
36		Dibenzo(a,h)anthracene	1.020	1.141	-11.9	201#	-0.01
37		Benzo(g,h,i)perylene	1.116	1.190	-6.6	195#	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 0